Partial Dynamical Symmetry in Nuclear Systems

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PARTIAL DYNAMICAL SYMMETRY IN NUCLEAR SYSTEMS

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Partial dynamical symmetry (PDS) extends and complements the concepts of exact and dynamical symmetry. It allows one to remove undesired constraints from an algebraic theory, while preserving some of the useful aspects of a dynamical symmetry, and to study the effects of symmetry breaking in a controlled manner. An example of a PDS in an interacting fermion system is presented. The associated PDS Hamiltonians are closely related with a realistic quadrupole-quadrupole interaction and provide new insights into this important interaction.

1. Introduction

Algebraic, symmetry-based theories provide an elegant and practical approach for describing and understanding a variety of physical systems. These theories offer the greatest simplifications when the interaction under consideration is symmetry preserving in the selected state labeling scheme, that is, when the Hamiltonian either commutes with all the generators of a particular group ('exact symmetry') or when it is written in terms of and commutes with the Casimir operators of a chain of nested groups ('dynamical symmetry'). An exact or dynamical symmetry not only facilitates the numerical treatment of the Hamiltonian, but also its interpretation and thus provides considerable insight into the physics of a given system.

Algebraic models can also be of value in situations where it is necessary to introduce symmetry-breaking terms in the Hamiltonian in order to achieve agreement with experimentally observed features. Pragmatically, it is often possible to decompose the offending terms into basic parts ("irreducible tensor operators") which exhibit specific transformation properties. Provided the appropriate group coupling coefficients and the matrix elements of some elementary tensor operators are available, matrix elements

of operators that connect inequivalent irreducible representations can be determined and the exact eigenvalues and eigenstates can then be obtained (at least in principle). Furthermore, by studying the effects of symmetry-breaking terms, one gains new insights into the dynamics of the system, the relevance and robustness of the symmetries under consideration, and their limitations. In some cases new symmetries emerge from a broken-symmetry regime. Their identification and interpretation is often simplified in the framework of an algebraic model. Examples of such emerging symmetries include quasi-dynamical symmetry¹ and pseudo-spin symmetry².

One can also consider introducing some intermediate structure that allows for symmetry breaking but retains some aspects of the dynamical symmetry. Partial dynamical symmetry (PDS) provides such a structure³. Two types of partial dynamical symmetry have been studied in recent years. Partial dynamical symmetry of the first kind preserves the advantages of a dynamical symmetry for a part of the system. It corresponds to a particular symmetry breaking for which the Hamiltonian is not invariant under the symmetry group and hence various irreducible representions (irreps) are mixed in its eigenstates, yet it possesses a subset of 'special' solvable states which respect the symmetry. PDS of the second kind corresponds to a symmetry breaking for which all eigenstates of the Hamiltonian preserve part of the dynamical symmetry⁴. In this scenario, the dynamical symmetry associated with an intermediate group G_2 in a subchain $G_1 \supset G_2 \supset G_3$ is broken for all states of the system, while the remaining (dynamical) symmetries are preserved. The resulting Hamiltonian is in general not analytically solvable, but its eigenstates can still be (partly) classified by quantum labels associated with the groups G_1 and G_3 . Further generalizations of the partial symmetry concept have been considered as well⁵.

This contribution will discuss an example of a partial dynamical symmetry in an interacting fermion system. More specifically, in the framework of the symplectic shell model (SSM), there exists a family of PDS Hamiltonians which are closely related to the nuclear quadrupole-quadrupole interaction. The Hamiltonians and their eigenstates are discussed and applications to the deformed light nuclei ²⁰Ne and ²⁴Mg are presented.

2. Quadrupole-Quadrupole Interaction in the SSM

The quadrupole-quadrupole interaction is an important ingredient in models that aim at reproducing quadrupole collective properties of nuclei. A model which is able to fully accommodate the action of the collec-

tive quadrupole operator, $Q_{2m} = \sqrt{\frac{16\pi}{5}} \sum_s r_s^2 Y_{2m}(\hat{r}_s)$, is the symplectic shell model (SSM), an algebraic scheme which respects the Pauli exclusion principle⁶. In the SSM, this operator takes the form $Q_{2m} = \sqrt{3}(\hat{C}_{2m}^{(11)} + \hat{A}_{2m}^{(20)} + \hat{B}_{2m}^{(02)})$, where $\hat{A}_{lm}^{(20)}$, $\hat{B}_{lm}^{(02)}$, and $\hat{C}_{lm}^{(11)}$ are symplectic generators with good SU(3) [superscript (λ, μ)] and SO(3) [subscript l, m] tensorial properties. The $\hat{A}_{lm}^{(20)}$ ($\hat{B}_{lm}^{(02)}$), l = 0 or 2, create (annihilate) $2\hbar\omega$ excitations in the system. The $\hat{C}_{lm}^{(11)}$, l = 1 or 2, generate a SU(3) subgroup and act only within one harmonic oscillator (h.o.) shell ($\sqrt{3}\hat{C}_{2m}^{(11)} = Q_{2m}^E$, the symmetrized quadrupole operator of Elliott, which does not couple different h.o. shells⁷, and $\hat{C}_{1m}^{(11)} = \hat{L}_m$, the orbital angular momentum operator). A fermion realization of these generators has been given⁸.

A basis for the symplectic model is generated by applying symmetrically coupled products of the $2\hbar\omega$ raising operator $\hat{A}^{(20)}$ with itself to the usual $0\hbar\omega$ many-particle shell-model states. Each $0\hbar\omega$ starting configuration is characterized by the distribution of oscillator quanta into the three cartesian directions, or, equivalently, by its $U(1)\times SU(3)$ quantum numbers $N_{\sigma}(\lambda_{\sigma}, \mu_{\sigma})$. Here $(\lambda_{\sigma}, \mu_{\sigma})$ are the Elliott SU(3) labels, and $N_{\sigma} = \sigma_1 + \sigma_2 + \sigma_3$ is related to the eigenvalue of the oscillator number operator. ²⁰Ne, for instance, has $N_{\sigma} = 48.5$ (after removal of the centerof-mass contribution) and $(\lambda_{\sigma}, \mu_{\sigma}) = (8,0)$. For ²⁴Mg, one finds $N_{\sigma} = 62.5$ and $(\lambda_{\sigma}, \mu_{\sigma}) = (8,4)$. The product of N/2 raising operators $\hat{A}^{(20)}$ generates $N\hbar\omega$ excitations for each starting irrep $N_{\sigma}(\lambda_{\sigma},\mu_{\sigma})$. Each such product operator $P^{N(\lambda_n,\mu_n)}$, labeled according to its SU(3) content, (λ_n,μ_n) , is coupled with $|N_{\sigma}(\lambda_{\sigma}, \mu_{\sigma})\rangle$ to good SU(3) symmetry $\rho(\lambda, \mu)$, with ρ denoting the multiplicity of the coupling $(\lambda_n, \mu_n) \otimes (\lambda_{\sigma}, \mu_{\sigma})$. To complete the basis state labeling, additional quantum numbers $\alpha = \kappa LM$ are required, where L denotes the angular momentum with projection M, and κ is a multiplicity index, which enumerates multiple occurrences of a particular L value in the SU(3) irrep (λ, μ) . The group chain corresponding to this labeling scheme is $Sp(6,R) \supset SU(3) \supset SO(3)$ which defines a dynamical symmetry basis.

The quadrupole-quadrupole interaction connects h.o. states differing in energy by $0\hbar\omega$, $\pm 2\hbar\omega$, and $\pm 4\hbar\omega$, and may be written as

$$Q_2 \cdot Q_2 = 9\hat{C}_{SU3} - 3\hat{C}_{Sp6} + \hat{H}_0^2 - 2\hat{H}_0 - 3\hat{L}^2 - 6\hat{A}_0\hat{B}_0 + \{\text{terms coupling different h.o. shells}\},$$
(1)

where \hat{C}_{SU3} and \hat{C}_{Sp6} are the quadratic Casimir invariants of SU(3) and Sp(6,R) with eigenvalues $2(\lambda^2 + \mu^2 + \lambda \mu + 3\lambda + 3\mu)/3$ and $2(\lambda^2_{\sigma} + \mu^2_{\sigma} + \lambda_{\sigma}\mu_{\sigma} + \mu^2_{\sigma} + \lambda_{\sigma}\mu_{\sigma})/3$

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 $3\lambda_{\sigma}+3\mu_{\sigma})/3+N_{\sigma}^2/3-4N_{\sigma}$, respectively. These operators, as well as the h.o. \hat{H}_0 and \hat{L}^2 terms, are diagonal in the dynamical symmetry basis. Unlike the Elliott quadrupole-quadrupole interaction, $Q_2^E \cdot Q_2^E = 6\hat{C}_{SU3} - 3\hat{L}^2$, the $Q_2 \cdot Q_2$ interaction of Eq. (1) breaks SU(3) symmetry within each h.o. shell since the term $\hat{A}_0\hat{B}_0 \equiv \hat{A}_0^{(20)}\hat{B}_0^{(02)} = (\{\hat{A}\times\hat{B}\}_0^{(00)} - \sqrt{5}\{\hat{A}\times\hat{B}\}_0^{(22)})/\sqrt{6}$ mixes different SU(3) irreps.

3. Partial Dynamical Symmetry in the SSM

In order to study the action of $Q_2 \cdot Q_2$ within a h.o. shell, we consider the following family of Hamiltonians:

$$H(\beta_0, \beta_2) = \beta_0 \hat{A}_0 \hat{B}_0 + \beta_2 \hat{A}_2 \cdot \hat{B}_2$$

$$= \frac{\beta_2}{18} (9\hat{C}_{SU3} - 9\hat{C}_{Sp6} + 3\hat{H}_0^2 - 36\hat{H}_0) + (\beta_0 - \beta_2)\hat{A}_0 \hat{B}_0 .$$
(2)

For $\beta_0 = \beta_2$, one recovers the dynamical symmetry, and with the special choice $\beta_0 = 12$, $\beta_2 = 18$, one obtains $Q_2 \cdot Q_2 = H(\beta_0 = 12, \beta_2 = 18) + const(N) - 3\hat{L}^2$ + terms coupling different shells, where const(N) is constant for a given h.o. $N\hbar\omega$ excitation.

It has been shown⁹ that $H(\beta_0, \beta_2)$ exhibits partial SU(3) symmetry of the first kind. Specifically, one finds that among the eigenstates of $H(\beta_0, \beta_2)$, there exists a subset of solvable pure-SU(3) states, the SU(3) \supset SO(3) classification of which depends on both the Elliott labels $(\lambda_{\sigma}, \mu_{\sigma})$ of the starting state and the symplectic excitation N. In general, one observes that all L-states in the starting configuration (N = 0) are solvable with good SU(3) symmetry $(\lambda_{\sigma}, \mu_{\sigma})$. For excited configurations (N > 0) and even one distinguishes between two possible cases:

- (a) $\lambda_{\sigma} > \mu_{\sigma}$: the pure states belong to $(\lambda, \mu) = (\lambda_{\sigma} N, \mu_{\sigma} + N)$ and have $L = \mu_{\sigma} + N, \mu_{\sigma} + N + 1, \dots, \lambda_{\sigma} N + 1$ with $N = 2, 4, \dots$ subject to $2N \leq (\lambda_{\sigma} \mu_{\sigma} + 1)$.
- (b) $\lambda_{\sigma} \leq \mu_{\sigma}$: the special states belong to $(\lambda, \mu) = (\lambda_{\sigma} + N, \mu_{\sigma})$ and have $L = \lambda_{\sigma} + N, \lambda_{\sigma} + N + 1, \dots, \lambda_{\sigma} + N + \mu_{\sigma}$ with $N = 2, 4, \dots$

The special states have well-defined symmetry $\operatorname{Sp}(6,\mathbb{R}) \supset \operatorname{SU}(3) \supset \operatorname{SO}(3)$ and are annihilated by \hat{B}_0 . This ensures that they are solvable eigenstates of $H(\beta_0, \beta_2)$ with eigenvalues E(N=0)=0, $E(N)=\beta_2 N(N_\sigma-\lambda_\sigma+\mu_\sigma-6+3N/2)/3$ for family (a), and $E(N)=\beta_2 N(N_\sigma+2\lambda_\sigma+\mu_\sigma-3+3N/2)/3$ for family (b). All $0\hbar\omega$ states are unmixed and span the entire $(\lambda_\sigma, \mu_\sigma)$ irrep. In contrast, for the excited levels (N>0), the pure states span only part of the corresponding $\operatorname{SU}(3)$ irreps. There are other states at each excited

level which do not preserve the SU(3) symmetry and therefore contain a mixture of SU(3) irreps. All eigenstates respect the Sp(6,R) and SO(3) symmetries. The partial SU(3) symmetry of $H(\beta_0, \beta_2)$ is converted into partial dynamical SU(3) symmetry by adding to it SO(3) rotation terms which lead to L(L+1)-type splitting but do not affect the wave functions. The solvable states form rotational bands and since their wave functions are known, one can evaluate the E2 rates between them⁹.

4. Applications

To illustrate that the PDS Hamiltonians discussed here are physically relevant, applications to realistic nuclear systems have been considered. Here the results for $^{20}{\rm Ne}$ and $^{24}{\rm Mg}$ are summarized. In particular, energy spectra and eigenstates of $H_{PDS}=h(N)+\xi H(\beta_0=12,\beta_2=18)+\gamma_2\hat{L}^2+\gamma_4\hat{L}^4$ are compared to those of $H_{Q\cdot Q}=\hat{H}_0-\chi Q_2\cdot Q_2+d_2\hat{L}^2+d_4\hat{L}^4,$ where h(N) is a constant for a given $N\hbar\omega$ excitation and contains the h.o. term \hat{H}_0 .

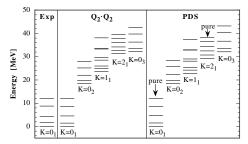


Figure 1. Energy spectra for 20 Ne. Comparison between experimental values (left), results from a symplectic $8\hbar\omega$ calculation (center) and a PDS calculation (right). The angular momenta of the positive parity states in the rotational bands are $L=0,2,4,\ldots$ for K=0 and $L=K,K+1,K+2,\ldots$ otherwise.

4.1. The ²⁰Ne Example

In Fig. 1, energy spectra of H_{PDS} are compared to those obtained from an $8\hbar\omega$ symplectic calculation (labeled $Q_2 \cdot Q_2$), and Fig. 2 shows the decomposition for representative (2^+) states of the five lowest rotational bands. The PDS Hamiltonian H_{PDS} acts only within one oscillator shell, hence its eigenfunctions do not contain admixtures from different $N\hbar\omega$ configurations. As expected, H_{PDS} has families of pure SU(3) eigenstates which can be organized into rotational bands, Fig. 1. The ground band belongs entirely to N=0, $(\lambda,\mu)=(8,0)$, and all states of the K=2₁ band have quantum labels N=2, $(\lambda,\mu)=(6,2)$, $\kappa=2$, see Fig. 2. A comparison

with the symplectic case shows that the $N\hbar\omega$ level to which a particular PDS band belongs is also dominant in the corresponding symplectic band. In addition, within this dominant excitation, eigenstates of H_{PDS} and $H_{Q,Q}$ have similar SU(3) distributions; in particular, both Hamiltonians favor the same $(\lambda, \mu)\kappa$ values. Significant differences in the structure of the wave functions appear, however, for the $K=0_2$ resonance band. In the symplectic calculation, this band contains almost equal contributions from the $0\hbar\omega$, $2\hbar\omega$, and $4\hbar\omega$ levels, with additional admixtures of $6\hbar\omega$ and $8\hbar\omega$ configurations, while in the PDS calculation, it belongs entirely to the $2\hbar\omega$ level. These structural differences are also evident in the interband transition $rates^9$ and reflect the action of the inter-shell coupling terms in Eq. (1). Increasing the strength χ of $Q_2 \cdot Q_2$ in $H_{Q \cdot Q}$ will also spread the other resonance bands over many $N\hbar\omega$ excitations. The K=2₁ band (which is pure in the PDS scheme) is found to resist this spreading more strongly than the other resonances. For physically relevant values of χ , the low-lying bands have the structure shown in Fig. 2.

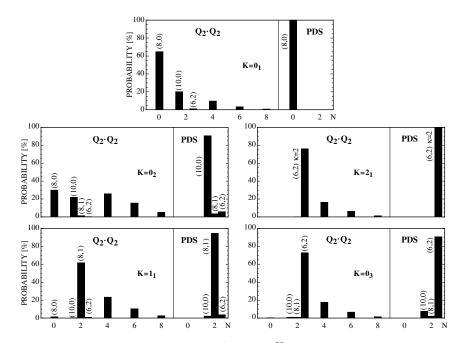


Figure 2. Decomposition for calculated 2^+ states of 20 Ne. Individual contributions from the relevant SU(3) irreps at the $0\hbar\omega$ and $2\hbar\omega$ levels are shown for both a symplectic $8\hbar\omega$ calculation (denoted $Q_2\cdot Q_2$) and a PDS calculation. In addition, the total strengths contributed by the $N\hbar\omega$ excitations for N>2 are given for the symplectic case.

4.2. The ²⁴Mg Example

For the triaxially deformed nucleus 24 Mg additional terms $\hat{X}_3 \equiv (\hat{L} \times Q^E) \cdot \hat{L}$ and $\hat{X}_4 \equiv (\hat{L} \times Q^E) \cdot (\hat{L} \times Q^E)$ are required in the Hamiltonian in order to reproduce the experimentally observed 'K-band splitting' between the ground and γ band of 24 Mg. Although these extra terms break the partial symmetry, for realistic interaction parameters the amount of symmetry breaking is very small ($\sim 1\%$). In Fig. 3, energy spectra of $H'_{PDS} = H_{PDS} + c_3 \hat{X}_3 + c_4 \hat{X}_4$ and $H'_{Q\cdot Q} = H_{Q\cdot Q} + c_3 \hat{X}_3 + c_4 \hat{X}_4$ are shown. H'_{PDS} has families of pure (and nearly pure) SU(3) eigenstates which can be organized into rotational bands; they are indicated in the figure.

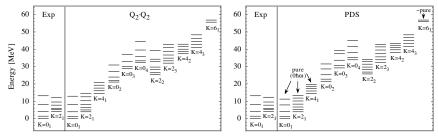


Figure 3. Energy spectra for 24 Mg. Energies from a PDS calculation (PDS) are compared to symplectic $6\hbar\omega$ results $(Q_2\cdot Q_2)$. Both $0\hbar\omega$ -dominated bands $(K=0_1,2_1,4_1)$ and some $2\hbar\omega$ resonance bands are shown. The $K=0_1$ and $K=2_1$ labels indicate the ground band and γ band, respectively.

The results are qualitatively similar to those for ²⁰Ne. The PDS Hamiltonian cannot account for intershell correlations, but it is able to reproduce various features of the quadrupole-quadrupole interaction, as can be seen in Fig. 4, where the structure of selected PDS eigenstates is compared to that of the corresponding $Q_2 \cdot Q_2$ eigenstates: PDS eigenfunctions do not contain admixtures from different $N\hbar\omega$ configurations, but belong entirely to one level of excitation. For reasonable interaction parameters, the $N\hbar\omega$ level to which a particular PDS band belongs is also dominant in the corresponding band of exact $Q_2 \cdot Q_2$ eigenstates. Within this dominant excitation, eigenstates of both Hamiltonians have similar SU(3) distributions. Structural differences, nevertheless, do arise and are reflected in the very sensitive interband transition rates⁹. Furthermore, due to the presence of \hat{X}_3 and \hat{X}_4, H'_{PDS} is only an approximate PDS Hamiltonian – the K=6₁ band has small admixtures from irreps other than $(\lambda, \mu) = (6, 6)$. Overall, it may be concluded that PDS eigenstates approximately reproduce the structure of the exact $Q_2 \cdot Q_2$ eigenstates, for both ground and resonance bands.

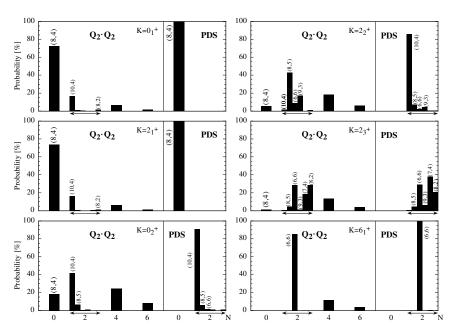


Figure 4. Decompositions for calculated $L^{\pi}=6^+$ states of ²⁴Mg. Eigenstates resulting from a symplectic $6\hbar\omega$ ($Q_2\cdot Q_2$) calculation are decomposed into their $0\hbar\omega$, $2\hbar\omega$, $4\hbar\omega$, $6\hbar\omega$ components. At the $0\hbar\omega$ and $2\hbar\omega$ levels, contributions from individual SU(3) irreps are shown, for higher excitations (N>2) only the summed strengths are given. Eigenstates of H'_{PDS} belong entirely to one $N\hbar\omega$ level of excitation, here $0\hbar\omega$ or $2\hbar\omega$; members of the K=0₁ and K=2₁ bands are pure and K=6₁ states are very nearly (>99%) pure.

5. Concluding Remarks

The notion of partial dynamical symmetry extends and complements the familiar concepts of exact and dynamical symmetry. It allows one to remove undesired constraints from an algebraic theory while preserving some of the useful aspects of a dynamical symmetry. As a result, the effects of symmetry breaking can be studied in a controlled manner and new insights into dynamics of the system under consideration are gained.

The work presented here focuses on a family of PDS Hamiltonians which are closely related to the deformation-inducing quadrupole-quadrupole interaction. For a particular parametrization, the PDS Hamiltonians take a form that is intermediate between the full quadrupole-quadrupole interaction, which couples states belonging to different harmonic oscillator shells, and the Elliott quadrupole-quadrupole interaction, which acts only within a shell. The intermediate scheme considered here extends the Elliott picture in that it includes (specific) SU(3) symmetry-breaking contributions.

At the same time, it is simpler than the full collective picture since it does not allow for mixing between different oscillator shells.

The PDS scheme sheds light on the in-shell behavior of the quadrupole-quadrupole interaction. For example, the symplectic model predicts the existence of states that are primarily dominated by one $N\hbar\omega$ level of excitation as well as states that contain strong multi-shell correlations^a. The states that resist the deformation-induced spreading over several $N\hbar\omega$ levels of excitation the strongest are those for which the associated PDS structure exhibits good (or almost good) SU(3) symmetry.

Acknowledgments

It has been a pleasure to present this work on the occasion of Jerry Draayer's 60th birthday. The research follows a tradition of using symmetry principles to gain insights into physical systems. Throughout his career, Jerry Draayer has contributed much to this tradition and has inspired many to seek the simplicity, symmetry, and beauty hidden within complex physical systems. The work presented here was carried out in collaboration with A. Leviatan (Hebrew University, Jerusalem). It was performed in part under the auspices of the U. S. Department of Energy by the University of California, Lawrence Livermore National Laboratory, under contract No. W-7405-Eng-48.

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^aThis result extends the findings of Bahri et~al. ¹⁰, who studied the structure of the giant quadrupole resonance in ²⁴Mg.